

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20

ring/chain bonds :

11-16

ring bonds :

1-2 1-6 1-12 2-3 2-14 3-4 4-5 5-6 5-7 6-11 7-8 8-9 9-10 10-11 12-13 13-14
15-17 15-16 16-20 17-18 18-19 19-20

exact/norm bonds :

1-2 1-6 1-12 2-3 2-14 3-4 4-5 5-6 5-7 6-11 7-8 8-9 9-10 10-11 11-16 12-13
13-14 15-17 15-16 16-20 17-18 18-19 19-20

G1:C,N

G2:C,O,S,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom
12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom

10/649,495

=> d his

(FILE 'HOME' ENTERED AT 11:52:45 ON 18 APR 2006)

FILE 'REGISTRY' ENTERED AT 11:54:28 ON 18 APR 2006

L1 STRUCTURE uploaded
L2 0 S L1
L3 26 S L1 SSS FUL

FILE 'CAPLUS' ENTERED AT 11:55:01 ON 18 APR 2006

L4 2 S L3
L5 1 S US20040058906/PN
SELECT RN L5 1-

FILE 'REGISTRY' ENTERED AT 11:55:35 ON 18 APR 2006

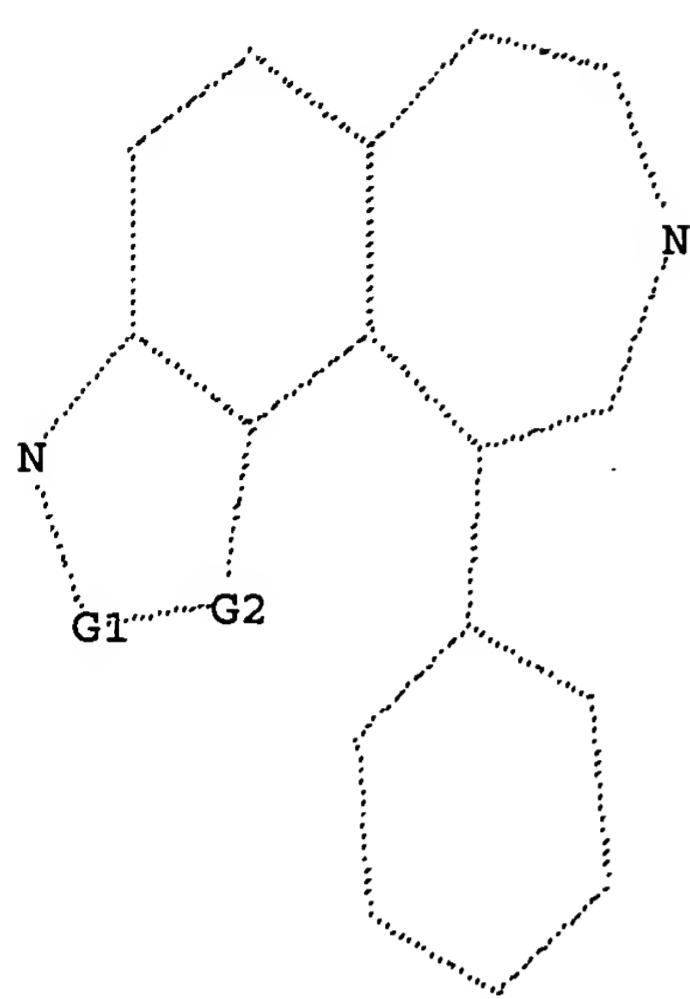
L6 76 S E1-76
L7 31 S L6 AND NRS=1
L8 29 S L6 AND NRS>1
L9 16 S L6 NOT (L7 OR L8)
L10 59174 S 5-6-7/SZ
L11 10 S L8 AND L10
L12 19 S L8 NOT L11
L13 1 S L12 AND 5-6-6-6-7/SZ
L14 18 S L12 NOT L13
L15 11 S L8 NOT L14
L16 12 S L7 AND 5-6-6-6-7/SZ
L17 19 S L7 NOT L16
L18 23 S L15 OR L16
L19 21 S L3 AND L18
L20 2 S L18 NOT L19
L21 STRUCTURE uploaded
L22 0 S L21
L23 29 S L21 SSS FUL

FILE 'CAPLUS' ENTERED AT 12:03:22 ON 18 APR 2006

L24 2 S L23

=> d l21

L21 HAS NO ANSWERS
L21 STR



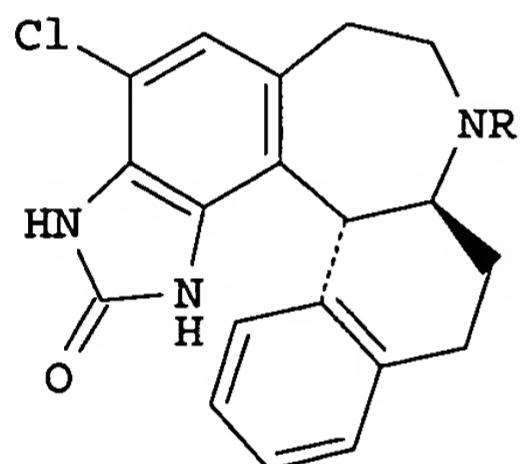
G1 C,N

G2 C,O,S,N

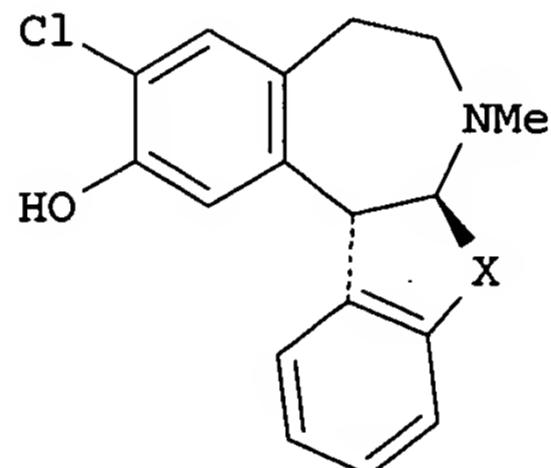
Structure attributes must be viewed using STN Express query preparation.

=> d ibib abs hitstr total

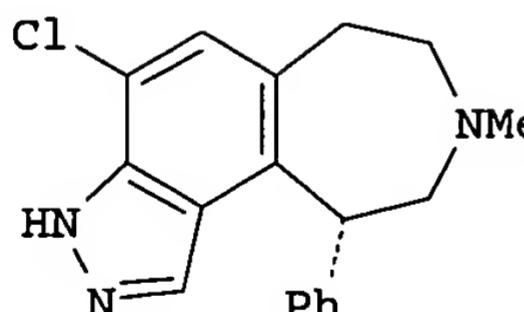
L24 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2006 ACS on STN
 ACCESSION NUMBER: 2005:31564 CAPLUS
 DOCUMENT NUMBER: 142:316682
 TITLE: Dopamine D1/D5 Receptor Antagonists with Improved Pharmacokinetics: Design, Synthesis, and Biological Evaluation of Phenol Bioisosteric Analogues of Benzazepine D1/D5 Antagonists
 AUTHOR(S): Wu, Wen-Lian; Burnett, Duane A.; Spring, Richard; Greenlee, William J.; Smith, Michelle; Favreau, Leonard; Fawzi, Ahmad; Zhang, Hongtao; Lachowicz, Jean E.
 CORPORATE SOURCE: Schering-Plough Research Institute, Kenilworth, NJ, 07033, USA
 SOURCE: Journal of Medicinal Chemistry (2005), 48(3), 680-693
 CODEN: JMCMAR; ISSN: 0022-2623
 PUBLISHER: American Chemical Society
 DOCUMENT TYPE: Journal
 LANGUAGE: English
 OTHER SOURCE(S): CASREACT 142:316682
 GI



I



II



III

AB Nonracemic fused benzazepines and naphthazepines such as I (R = H, Me) are prepared as selective dopamine D1 and D5 receptor antagonists with improved bioavailability over related high affinity dopamine D1 and D5 receptor antagonists by replacement of the phenol moiety in II (X = CH₂CH₂) with a variety of fused hydrogen-bond donating moieties. Benzazepines in which the hydrogen bond donor is pointed approx. parallel to an axis through the benzazepine nitrogen and the benzo ring are more effective as selective dopamine D1 and D5 receptor antagonists than benzazepines in which the hydrogen bond donor is pointed away from the axis. Attempts to replace the phenol group in a benzazepine II (X = H₂) with a bioisostere lead to decreased binding to the desired dopamine receptors; an indazolobenzazepine III is an active dopamine D1 and D5 receptor antagonist. I (R = H, Me) show improved pharmacokinetic behavior over II (X = CH₂CH₂) in rats; III shows similar pharmacokinetic behavior in rats

to II (X = H2).

IT 668476-37-7P 668476-46-8P

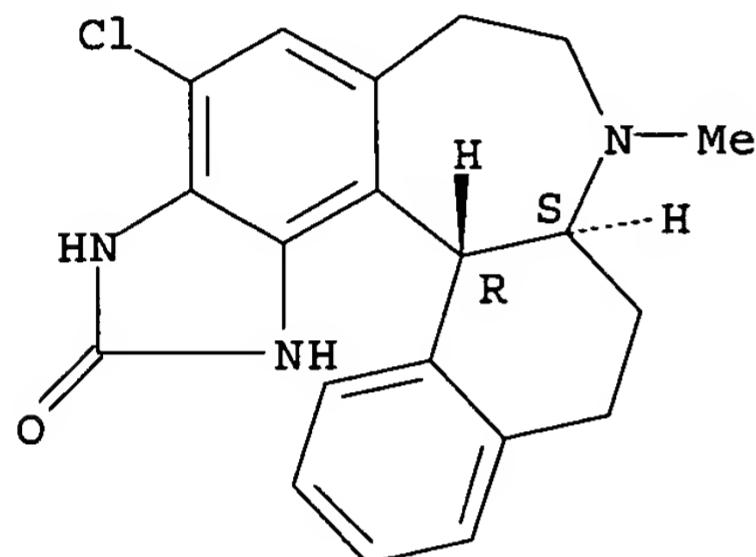
RL: PAC (Pharmacological activity); PKT (Pharmacokinetics); RCT (Reactant); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent)

(preparation of nonracemic benzazepines and naphthazepines containing bioisosteric replacements for a phenol moiety and their activity as selective dopamine D1/D5 receptor antagonists and the pharmacokinetic behavior of selected benzazepines)

RN 668476-37-7 CAPLUS

CN Benzimidazo[5,4-d]benzo[g][1]benzazepin-2(1H)-one, 4-chloro-3,6,7,8,8a,9,10,14b-octahydro-8-methyl-, (8aS,14bR)- (9CI) (CA INDEX NAME)

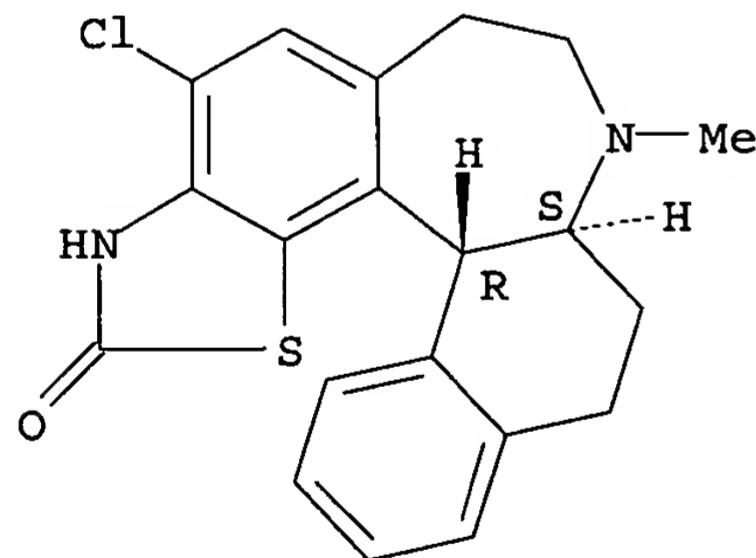
Absolute stereochemistry.



RN 668476-46-8 CAPLUS

CN 2H-Benzo[g]benzothiazolo[6,7-d][1]benzazepin-2-one, 4-chloro-3,6,7,8,8a,9,10,14b-octahydro-8-methyl-, (8aS,14bR)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



IT 668476-25-3P 668476-47-9P 847265-03-6P

RL: PAC (Pharmacological activity); PKT (Pharmacokinetics); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)

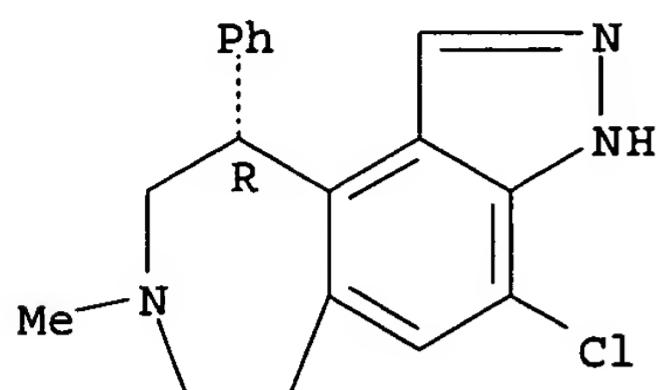
(preparation of nonracemic benzazepines and naphthazepines containing bioisosteric replacements for a phenol moiety and their activity as selective dopamine D1/D5 receptor antagonists and the pharmacokinetic behavior of selected benzazepines)

RN 668476-25-3 CAPLUS

CN Pyrazolo[4,3-g][3]benzazepine, 4-chloro-3,6,7,8,9,10-hexahydro-8-methyl-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

10/649,495

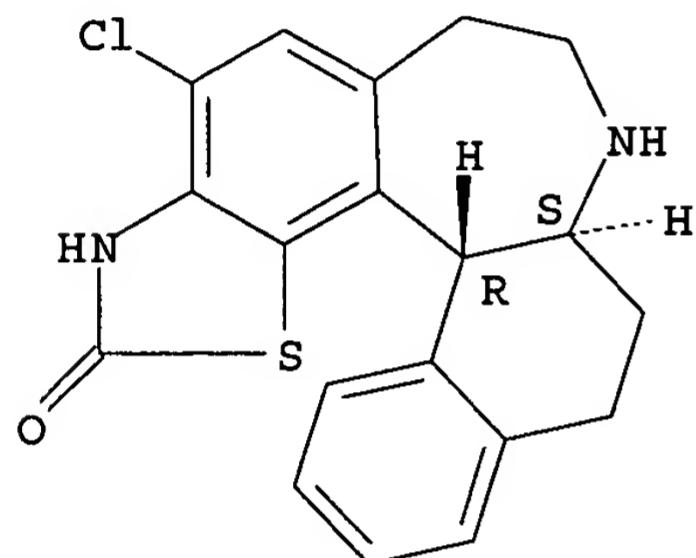
Absolute stereochemistry.



RN 668476-47-9 CAPLUS

CN 2H-Benz[*g*]benzothiazolo[6,7-*d*][1]benzazepin-2-one, 4-chloro-3,6,7,8,8a,9,10,14b-octahydro-, (8a*S*,14b*R*)- (9CI) (CA INDEX NAME)

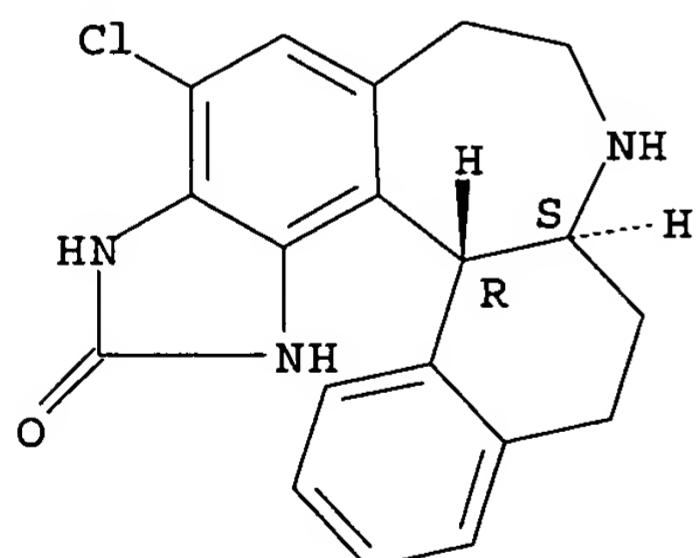
Absolute stereochemistry.



RN 847265-03-6 CAPLUS

CN Benzimidazo[5,4-d]benzo[g][1]benzazepin-2(1H)-one, 4-chloro-3,6,7,8,8a,9,10,14b-octahydro-, monohydrochloride, (8a*S*,14b*R*)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



● HCl

IT 668476-24-2P 668476-44-6P 668476-57-1P

668476-66-2P

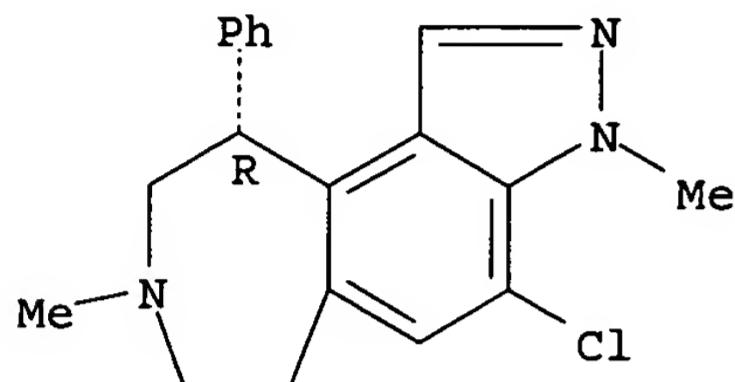
RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent)

(preparation of nonracemic benzazepines and naphthazepines containing bioisosteric replacements for a phenol moiety and their activity as selective dopamine D₁/D₅ receptor antagonists and the pharmacokinetic behavior of selected benzazepines)

RN 668476-24-2 CAPLUS

CN Pyrazolo[4,3-g] [3]benzazepine, 4-chloro-3,6,7,8,9,10-hexahydro-3,8-dimethyl-10-phenyl-, (10R) - (9CI) (CA INDEX NAME)

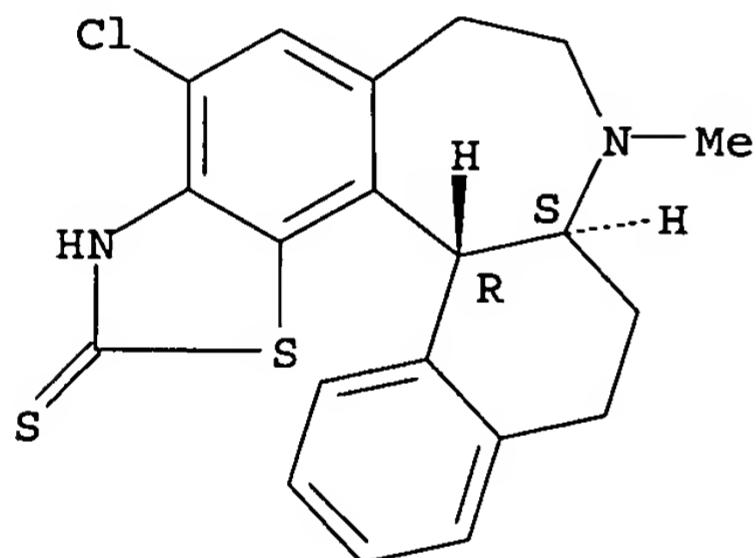
Absolute stereochemistry.



RN 668476-44-6 CAPLUS

CN 2H-Benzo[g]benzothiazolo[6,7-d] [1]benzazepine-2-thione, 4-chloro-3,6,7,8,8a,9,10,14b-octahydro-8-methyl-, (8aS,14bR) - (9CI) (CA INDEX NAME)

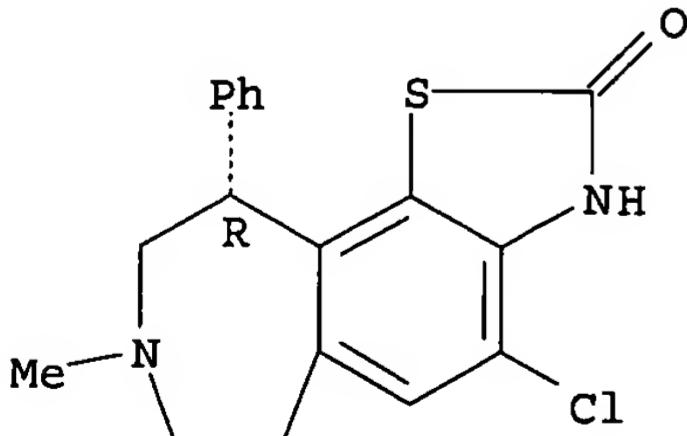
Absolute stereochemistry.



RN 668476-57-1 CAPLUS

CN 2H-Thiazolo[5,4-g] [3]benzazepin-2-one, 4-chloro-3,6,7,8,9,10-hexahydro-8-methyl-10-phenyl-, (10R) - (9CI) (CA INDEX NAME)

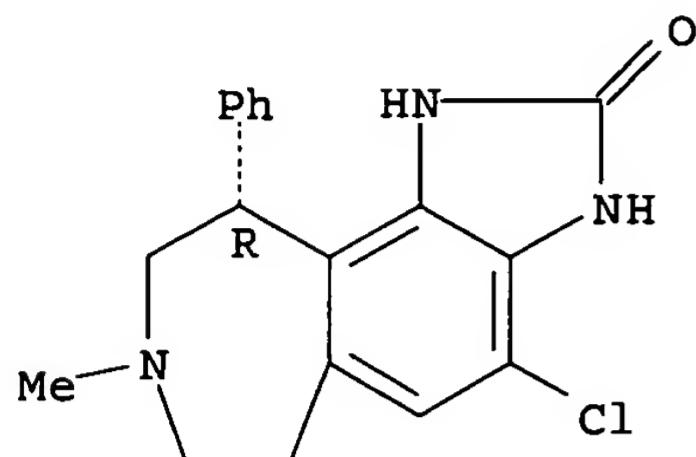
Absolute stereochemistry.



RN 668476-66-2 CAPLUS

CN Imidazo[4,5-g] [3]benzazepin-2(1H)-one, 4-chloro-3,6,7,8,9,10-hexahydro-8-methyl-10-phenyl-, (10R) - (9CI) (CA INDEX NAME)

Absolute stereochemistry.



IT 668476-26-4P 668476-29-7P 668476-34-4P
 668476-36-6P 668476-38-8P 668476-58-2P
 668476-67-3P 668476-68-4P

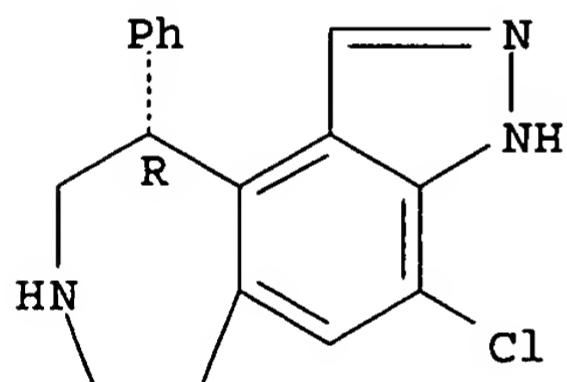
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)

(preparation of nonracemic benzazepines and naphthazepines containing bioisosteric replacements for a phenol moiety and their activity as selective dopamine D1/D5 receptor antagonists and the pharmacokinetic behavior of selected benzazepines)

RN 668476-26-4 CAPLUS

CN Pyrazolo[4,3-g][3]benzazepine, 4-chloro-3,6,7,8,9,10-hexahydro-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

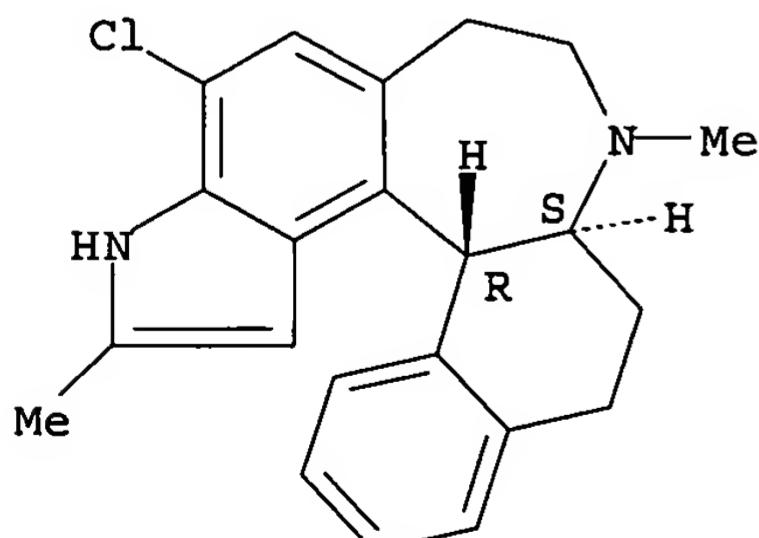
Absolute stereochemistry.



RN 668476-29-7 CAPLUS

CN Benz[g]indolo[5,4-d][1]benzazepine, 7-chloro-1,2,2a,3,4,5,8,10c-octahydro-3,9-dimethyl-, (2aS,10cR)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

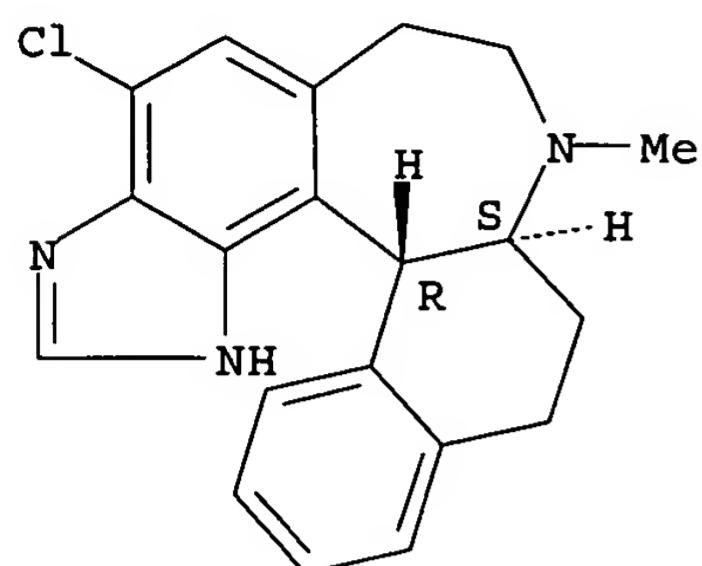


RN 668476-34-4 CAPLUS

CN Benzimidazo[5,4-d]benzo[g][1]benzazepine, 4-chloro-1,6,7,8,8a,9,10,14b-

octahydro-8-methyl-, (8aS,14bR)- (9CI) (CA INDEX NAME)

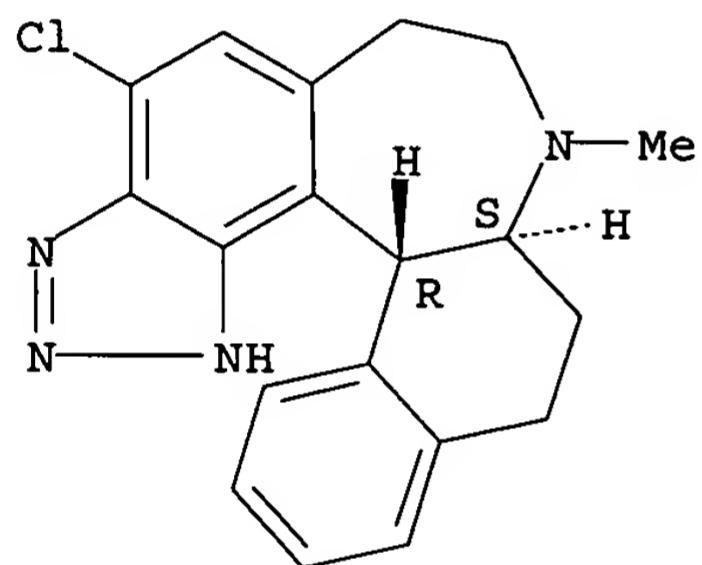
Absolute stereochemistry.



RN 668476-36-6 CAPLUS

CN Benzo [g] benzotriazolo [5,4-d] [1]benzazepine, 4-chloro-1,6,7,8,8a,9,10,14b-octahydro-8-methyl-, (8aS,14bR)- (9CI) (CA INDEX NAME)

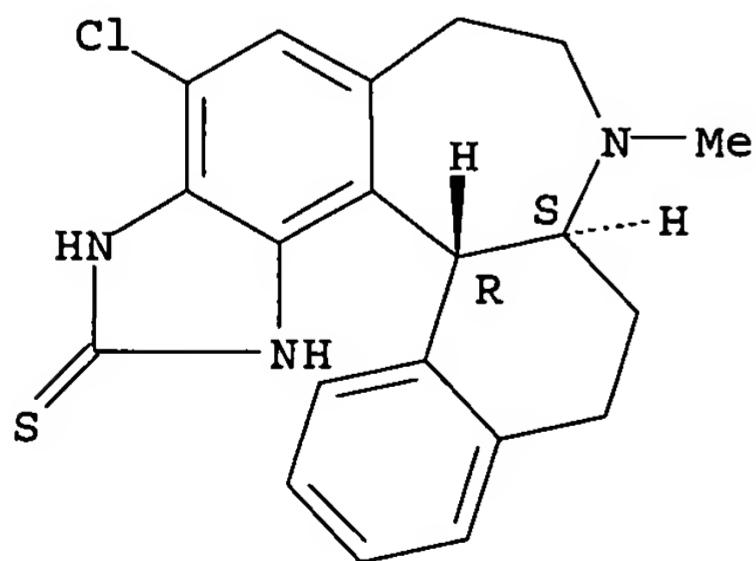
Absolute stereochemistry.



RN 668476-38-8 CAPLUS

CN Benzimidazo [5,4-d]benzo [g] [1]benzazepine-2(1H)-thione, 4-chloro-3,6,7,8,8a,9,10,14b-octahydro-8-methyl-, (8aS,14bR)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

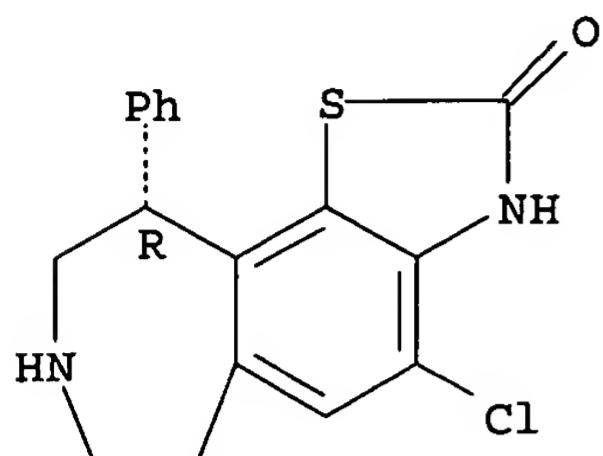


RN 668476-58-2 CAPLUS

CN 2H-Thiazolo[5,4-g][3]benzazepin-2-one, 4-chloro-3,6,7,8,9,10-hexahydro-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

10/649,495

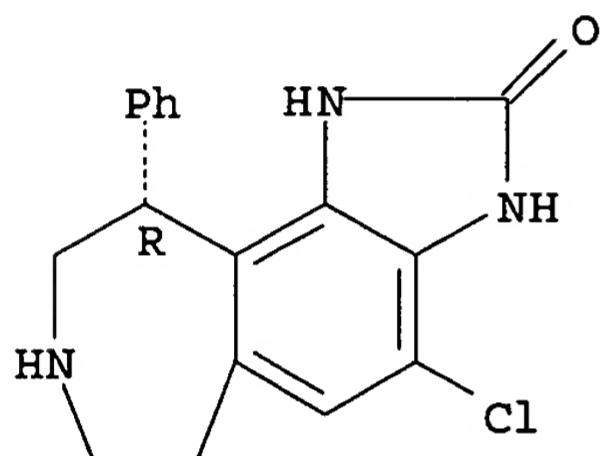
Absolute stereochemistry.



RN 668476-67-3 CAPLUS

CN Imidazo[4,5-g][3]benzazepin-2(1H)-one, 4-chloro-3,6,7,8,9,10-hexahydro-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

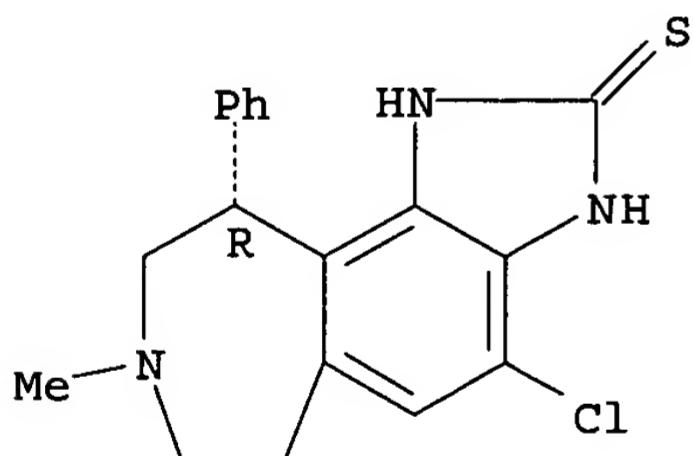
Absolute stereochemistry.



RN 668476-68-4 CAPLUS

CN Imidazo[4,5-g][3]benzazepine-2(1H)-thione, 4-chloro-3,6,7,8,9,10-hexahydro-8-methyl-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



IT 847265-05-8

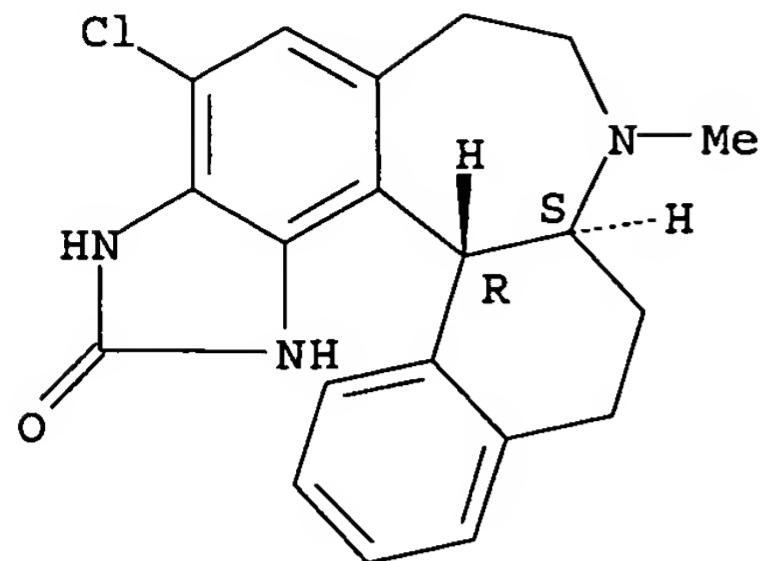
RL: RCT (Reactant); RACT (Reactant or reagent)

(preparation of nonracemic benzazepines and naphthazepines containing bioisosteric replacements for a phenol moiety and their activity as selective dopamine D1/D5 receptor antagonists and the pharmacokinetic behavior of selected benzazepines)

RN 847265-05-8 CAPLUS

CN Benzimidazo[5,4-d]benzo[g][1]benzazepin-2(1H)-one, 4-chloro-3,6,7,8,8a,9,10,14b-octahydro-8-methyl-, monohydrochloride, (8aS,14bR)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



● HCl

IT 668476-45-7P 668476-55-9P 668476-56-0P

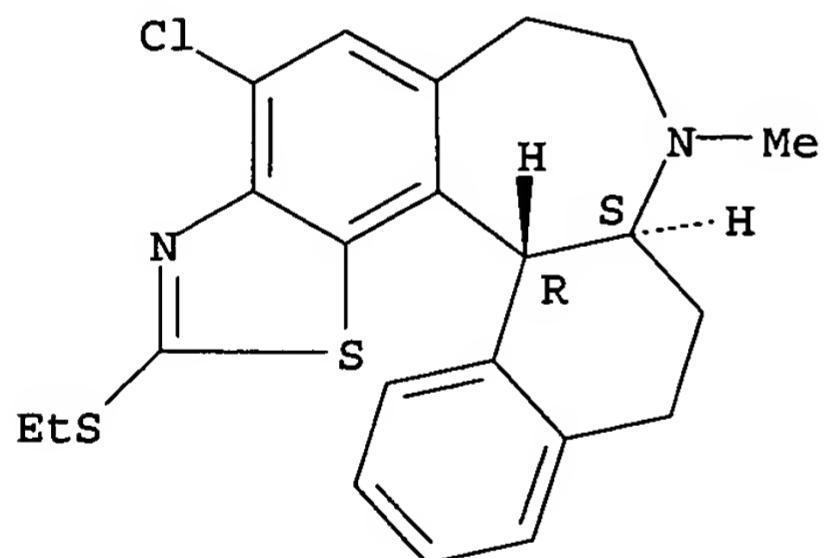
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of nonracemic benzazepines and naphthazepines containing bioisosteric replacements for a phenol moiety and their activity as selective dopamine D₁/D₅ receptor antagonists and the pharmacokinetic behavior of selected benzazepines)

RN 668476-45-7 CAPLUS

CN 6H-Benzo[g]benzothiazolo[6,7-d][1]benzazepine, 4-chloro-2-(ethylthio)-7,8,8a,9,10,14b-hexahydro-8-methyl-, (8aS,14bR)- (9CI) (CA INDEX NAME)

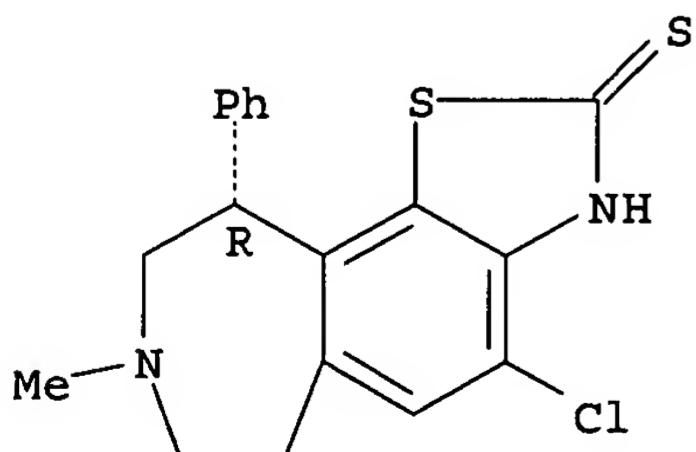
Absolute stereochemistry.



RN 668476-55-9 CAPLUS

CN 2H-Thiazolo[5,4-g][3]benzazepine-2-thione, 4-chloro-3,6,7,8,9,10-hexahydro-8-methyl-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

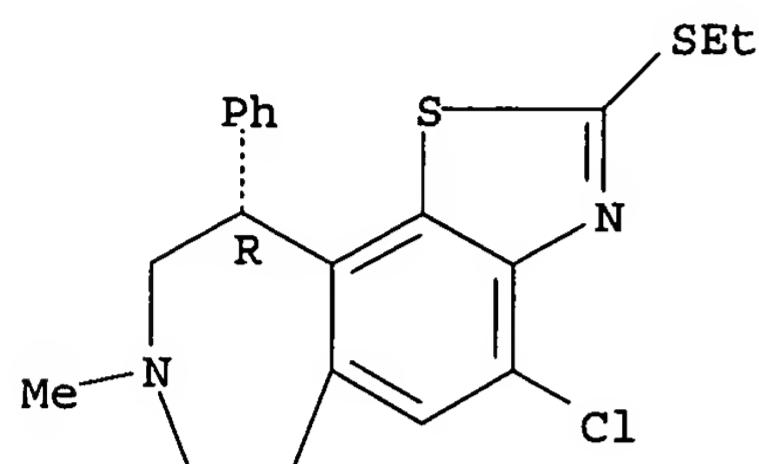


10/649,495

RN 668476-56-0 CAPLUS

CN 6H-Thiazolo[5,4-g][3]benzazepine, 4-chloro-2-(ethylthio)-7,8,9,10-tetrahydro-8-methyl-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



REFERENCE COUNT:

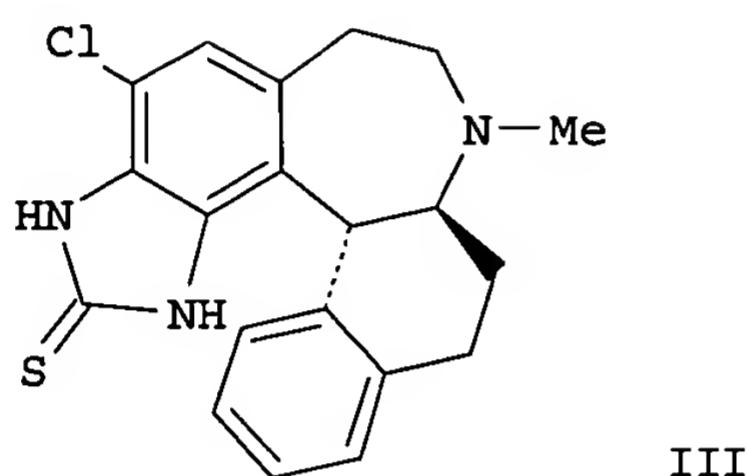
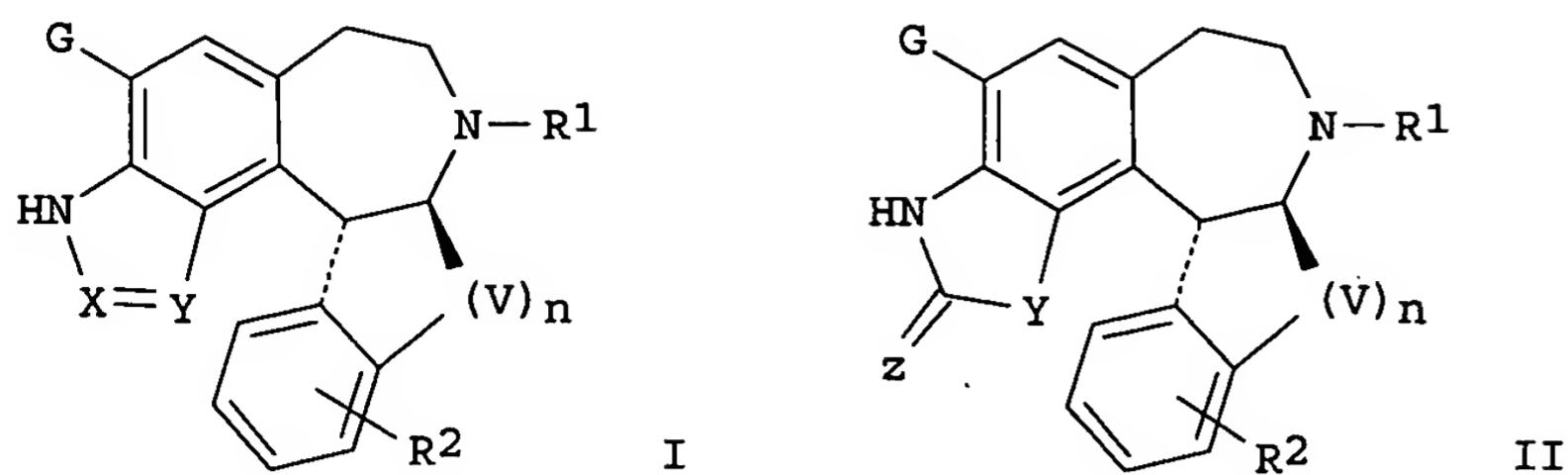
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THERE ARE 60 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

10/649,495

L24 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2006 ACS on STN
ACCESSION NUMBER: 2004:203836 CAPLUS
DOCUMENT NUMBER: 140:235618
TITLE: Preparation of benzazepine derivatives as selective D1/D5 receptor antagonists for the treatment of obesity and CNS disorders
INVENTOR(S): Wu, Wen-Lian; Burnett, Duane A.; Greenlee, William J.; Sasikumar, Thavalakulam K.
PATENT ASSIGNEE(S): Schering Corporation, USA
SOURCE: PCT Int. Appl., 54 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

| PATENT NO. | KIND | DATE | APPLICATION NO. | DATE |
|---|------|-------------------|-----------------|------------|
| WO 2004020442 | A1 | 20040311 | WO 2003-US26878 | 20030827 |
| W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, HR, HU, ID, IL, IN, IS, JP, KG, KR, KZ, LC, LK, LR, LT, LU, LV, MA, MD, MG, MK, MN, MX, MZ, NI, NO, NZ, PG, PH, PL, PT, RO, RU, SC, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UZ, VC, VN, YU, ZA, ZM | | | | |
| RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG | | | | |
| CA 2495879 | AA | 20040311 | CA 2003-2495879 | 20030827 |
| AU 2003262926 | A1 | 20040319 | AU 2003-262926 | 20030827 |
| US 2004058906 | A1 | 20040325 | US 2003-649495 | 20030827 |
| EP 1537115 | A1 | 20050608 | EP 2003-791865 | 20030827 |
| R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK | | | | |
| JP 2006501244 | T2 | 20060112 | JP 2004-531584 | 20030827 |
| PRIORITY APPLN. INFO.: | | | US 2002-406856P | P 20020829 |
| | | | WO 2003-US26878 | W 20030827 |
| OTHER SOURCE(S): GI | | MARPAT 140:235618 | | |



AB The title compds. I and II [R1 = H, alkyl, allyl, cycloalkyl, cycloalkyl(alkyl); R2 = H, halo, alkyl, alkylthio, alkylsulfonyl, OH, alkoxy, CF₃, CF₃O, aryl, CHO, NO₂, substituted amines, CN, etc.; G = H, halo, alkyl, alkylthio, NO₂, CN, OH, alkoxy, alkylsulfinyl, alkylsulfonyl, CF₃, or CF₃O; X = CH, C(alkyl), CCF₃, or N; Y = CH, C(alkyl) or N; Z = NH, N(alkyl), S, or O; V = CH₂; n = 0-2, when n = 0, the carbons to which (V)_n is shown connected are not linked to each other but are linked to hydrogen] were prepared as selective D₁/D₅ receptor antagonists for the treatment of obesity and CNS disorders. For example, treatment of (6aS,13bR)-12,13-diamino-11-chloro-6,6a,7,8,9,13b-hexahydro-7-methyl-5H-benzo[d]naphth[2,1-b]azepine (preparation given) with thiocarbonyldiimidazole gave compound III. The latter is a novel antagonist for D₁ receptor with K_i = 40 nM.

IT 668476-46-8P 668476-57-1P 668476-64-0P

668476-66-2P

RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP

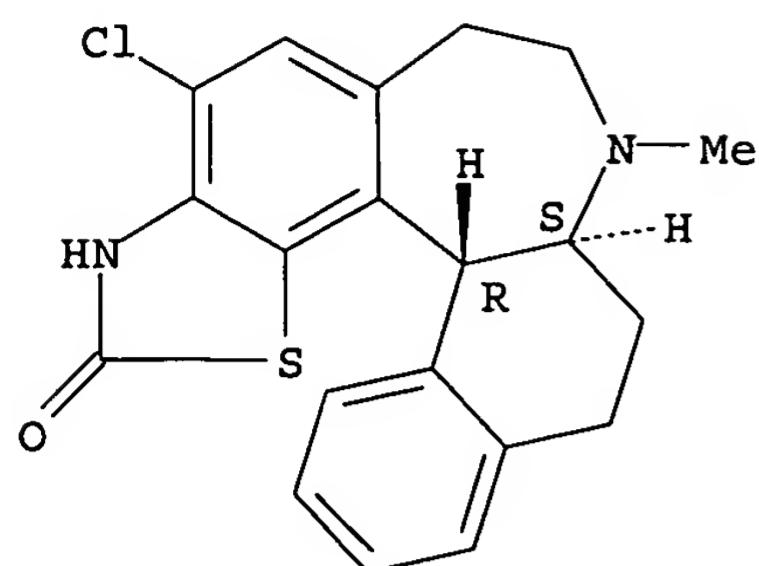
Preparation); RACT (Reactant or reagent); USES (Uses)

(preparation of benzazepine derivs. as selective D1/D5 receptor antagonists for the treatment of obesity and CNS disorders)

RN 668476-46-8 CAPLUS

CN 2H-Benzo[g]benzothiazolo[6,7-d][1]benzazepin-2-one, 4-chloro-
3,6,7,8,8a,9,10,14b-octahydro-8-methyl-, (8aS,14bR)- (9CI) (CA INDEX
NAME)

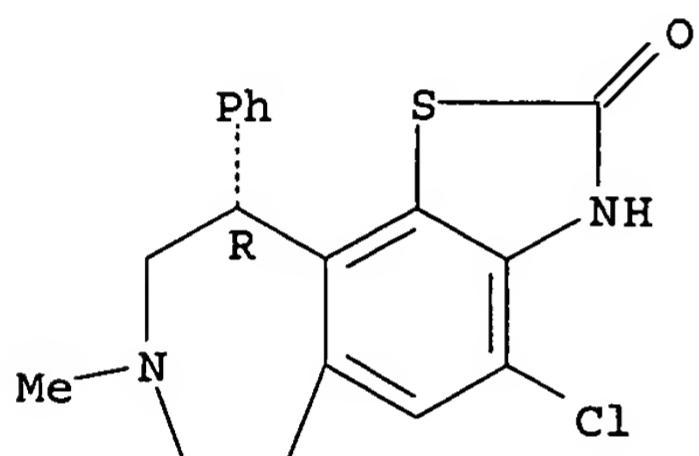
Absolute stereochemistry.



RN 668476-57-1 CAPLUS

CN 2H-Thiazolo[5,4-g][3]benzazepin-2-one, 4-chloro-3,6,7,8,9,10-hexahydro-8-methyl-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

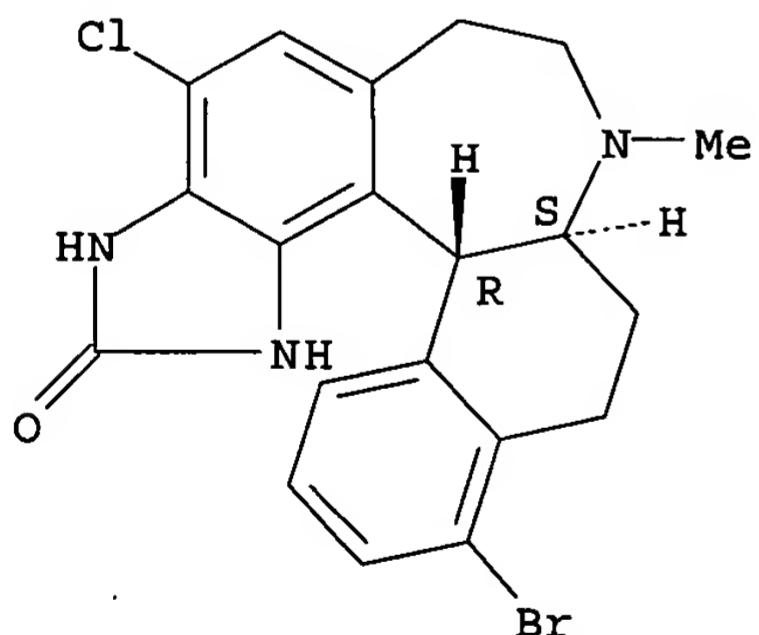
Absolute stereochemistry.



RN 668476-64-0 CAPLUS

CN Benzimidazo[5,4-d]benzo[g][1]benzazepin-2(1H)-one, 11-bromo-4-chloro-3,6,7,8,8a,9,10,14b-octahydro-8-methyl-, (8aS,14bR)- (9CI) (CA INDEX NAME)

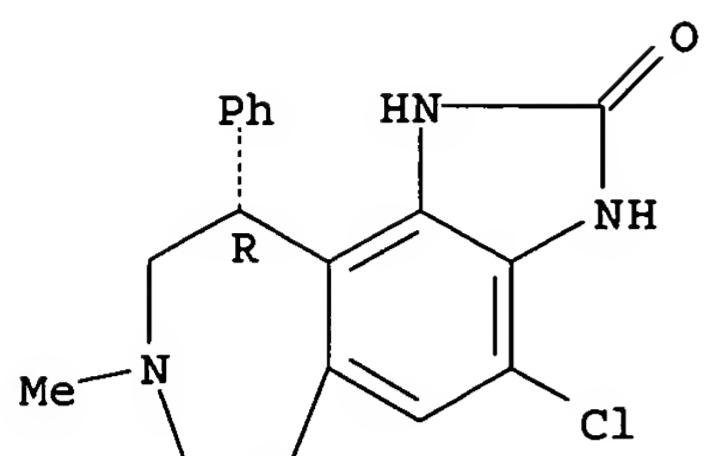
Absolute stereochemistry.



RN 668476-66-2 CAPLUS

CN Imidazo[4,5-g][3]benzazepin-2(1H)-one, 4-chloro-3,6,7,8,9,10-hexahydro-8-methyl-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



IT 668476-25-3P 668476-26-4P 668476-29-7P
 668476-34-4P 668476-36-6P 668476-38-8P
 668476-39-9P 668476-47-9P 668476-48-0P
 668476-58-2P 668476-65-1P 668476-67-3P
 668476-68-4P

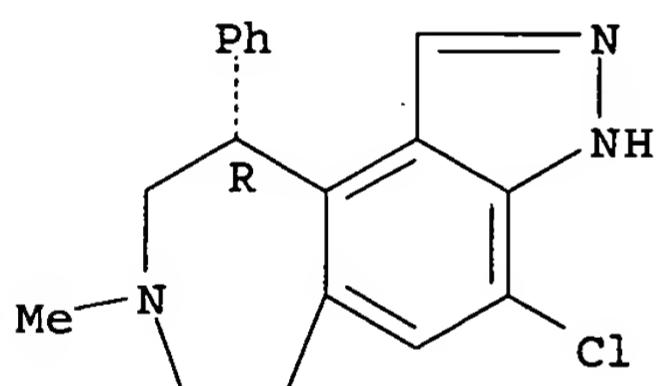
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of benzazepine derivs. as selective D1/D5 receptor antagonists for the treatment of obesity and CNS disorders)

RN 668476-25-3 CAPLUS

CN Pyrazolo[4,3-g] [3]benzazepine, 4-chloro-3,6,7,8,9,10-hexahydro-8-methyl-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

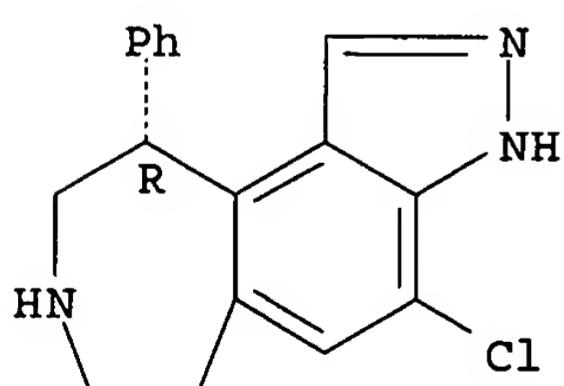
Absolute stereochemistry.



RN 668476-26-4 CAPLUS

CN Pyrazolo[4,3-g] [3]benzazepine, 4-chloro-3,6,7,8,9,10-hexahydro-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

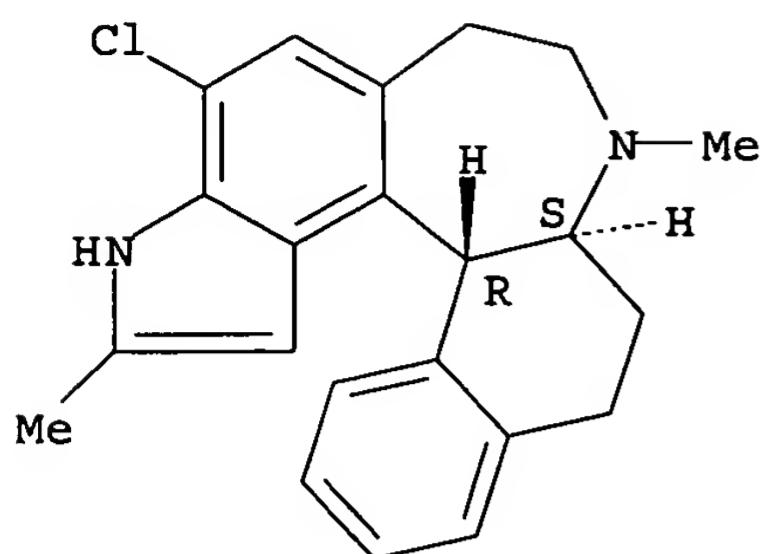
Absolute stereochemistry.



RN 668476-29-7 CAPLUS

CN Benz[g]indolo[5,4-d] [1]benzazepine, 7-chloro-1,2,2a,3,4,5,8,10c-octahydro-3,9-dimethyl-, (2aS,10cR)- (9CI) (CA INDEX NAME)

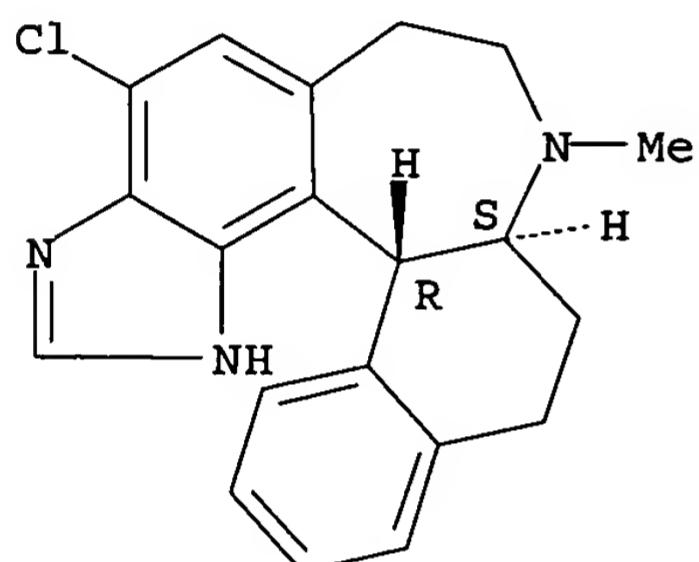
Absolute stereochemistry.



RN 668476-34-4 CAPLUS

CN Benzimidazo[5,4-d]benzo[g][1]benzazepine, 4-chloro-1,6,7,8,8a,9,10,14b-octahydro-8-methyl-, (8aS,14bR)- (9CI) (CA INDEX NAME)

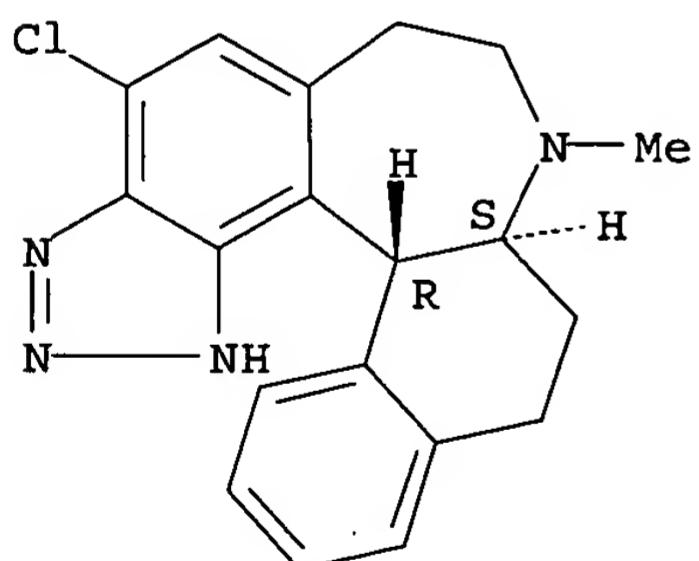
Absolute stereochemistry.



RN 668476-36-6 CAPLUS

CN Benzo[g]benzotriazolo[5,4-d] [1]benzazepine, 4-chloro-1,6,7,8,8a,9,10,14b-octahydro-8-methyl-, (8aS,14bR) - (9CI) (CA INDEX NAME)

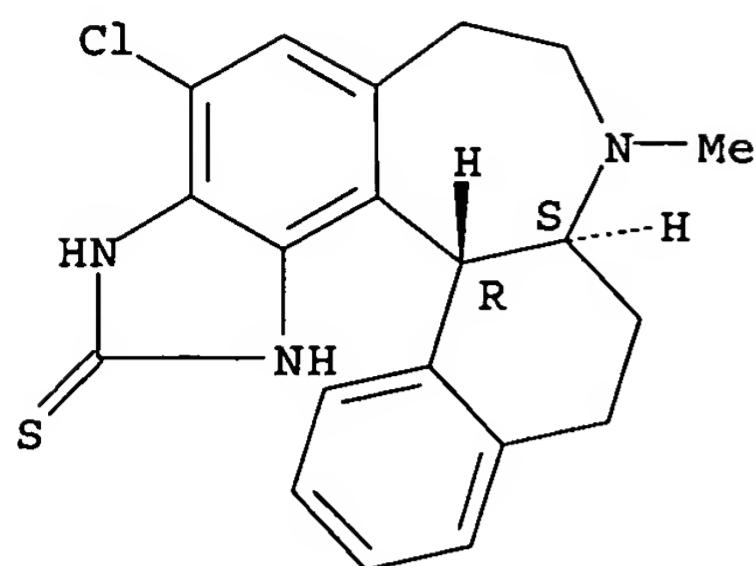
Absolute stereochemistry.



RN 668476-38-8 CAPLUS

CN Benzimidazo[5,4-d]benzo[g][1]benzazepine-2(1H)-thione,
4-chloro-3,6,7,8,8a,9,10,14b-octahydro-8-methyl-, (8aS,14bR)- (9CI) (CA
INDEX NAME)

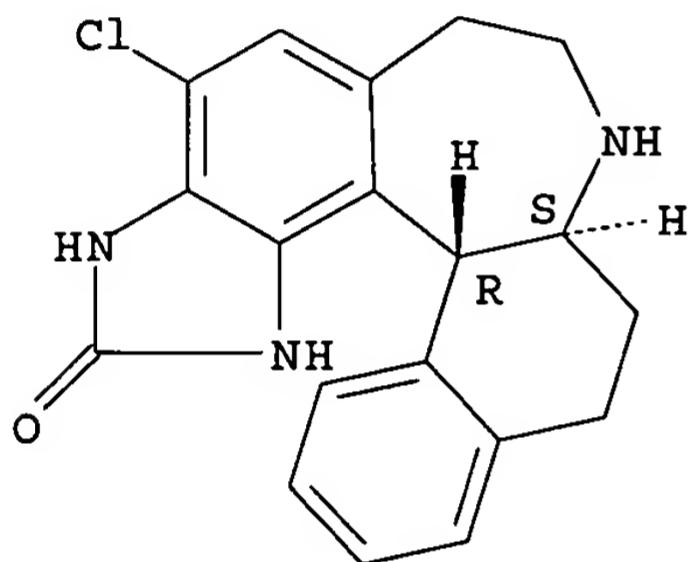
Absolute stereochemistry.



RN 668476-39-9 CAPLUS

CN Benzimidazo[5,4-d]benzo[g][1]benzazepin-2(1H)-one, 4-chloro-3,6,7,8,8a,9,10,14b-octahydro-, (8aS,14bR)- (9CI) (CA INDEX NAME)

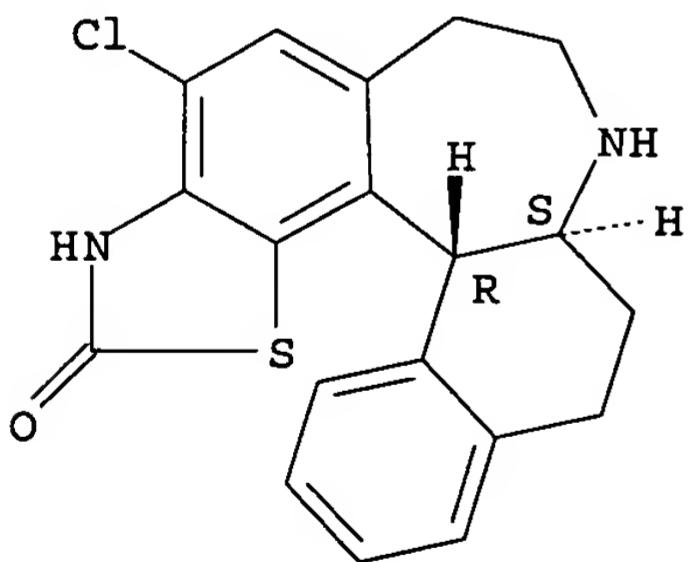
Absolute stereochemistry.



RN 668476-47-9 CAPLUS

CN 2H-Benzothiazolo[6,7-d][1]benzazepin-2-one, 4-chloro-3,6,7,8,8a,9,10,14b-octahydro-, (8aS,14bR)- (9CI) (CA INDEX NAME)

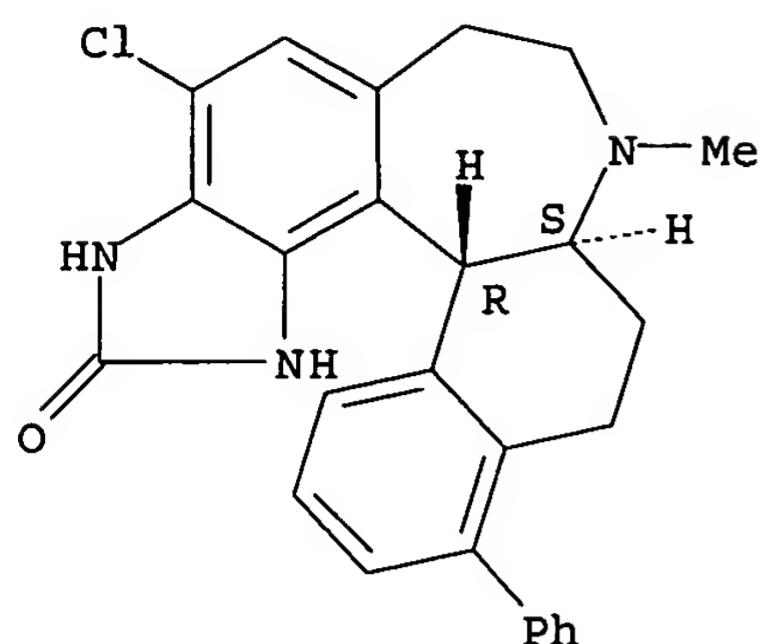
Absolute stereochemistry.



RN 668476-48-0 CAPLUS

CN Benzimidazo[5,4-d]benzo[g][1]benzazepin-2(1H)-one, 4-chloro-3,6,7,8,8a,9,10,14b-octahydro-8-methyl-11-phenyl-, (8aS,14bR)- (9CI) (CA INDEX NAME)

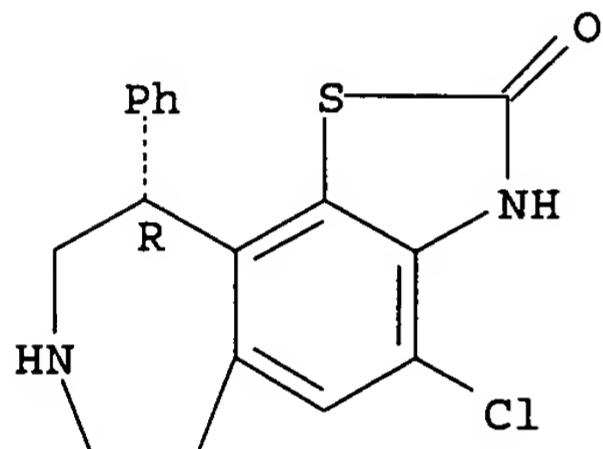
Absolute stereochemistry.



RN 668476-58-2 CAPLUS

CN 2H-Thiazolo[5,4-g][3]benzazepin-2-one, 4-chloro-3,6,7,8,9,10-hexahydro-10-phenyl-, (10R) - (9CI) (CA INDEX NAME)

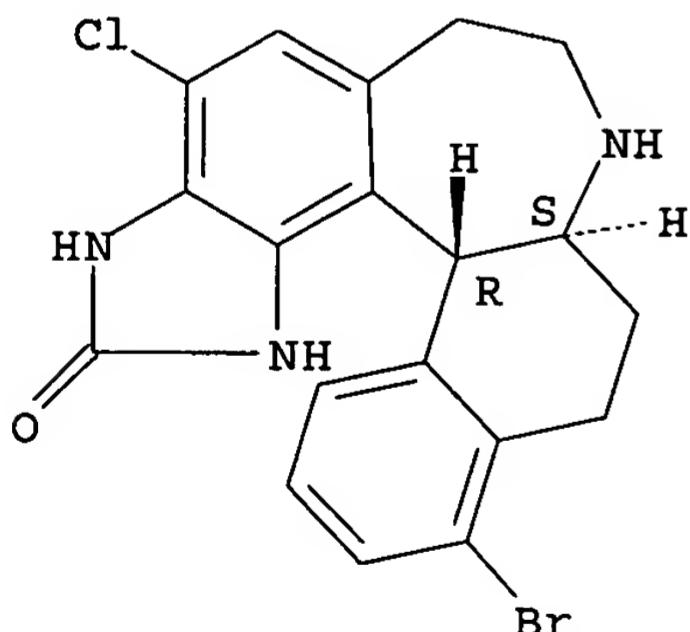
Absolute stereochemistry.



RN 668476-65-1 CAPLUS

CN Benzimidazo[5,4-d]benzo[g][1]benzazepin-2(1H)-one, 11-bromo-4-chloro-3,6,7,8,8a,9,10,14b-octahydro-, (8aS,14bR) - (9CI) (CA INDEX NAME)

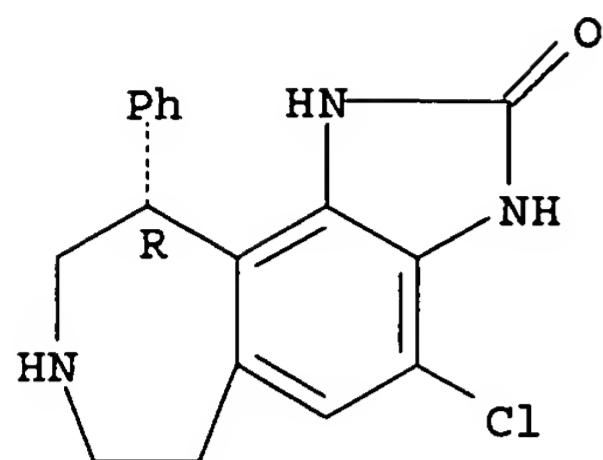
Absolute stereochemistry.



RN 668476-67-3 CAPLUS

CN Imidazo[4,5-g][3]benzazepin-2(1H)-one, 4-chloro-3,6,7,8,9,10-hexahydro-10-phenyl-, (10R) - (9CI) (CA INDEX NAME)

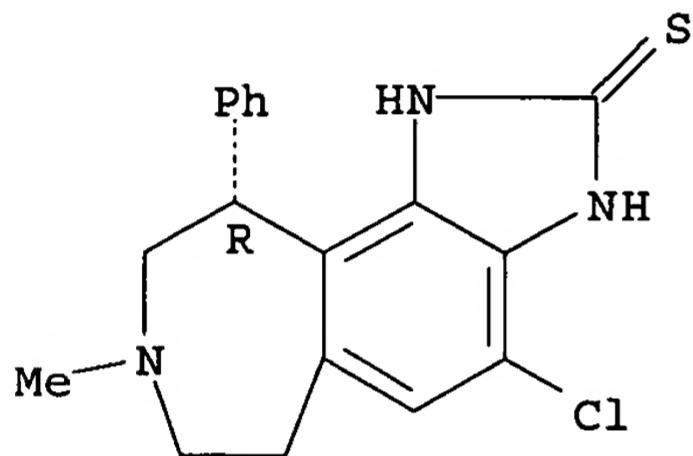
Absolute stereochemistry.



RN 668476-68-4 CAPLUS

CN Imidazo[4,5-g][3]benzazepine-2(1H)-thione, 4-chloro-3,6,7,8,9,10-hexahydro-8-methyl-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 668476-24-2P 668476-37-7P 668476-44-6P
668476-45-7P 668476-55-9P 668476-56-0P

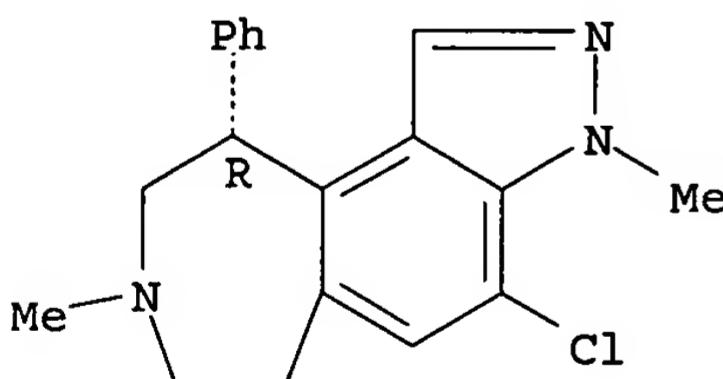
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of benzazepine derivs. as selective D1/D5 receptor antagonists for the treatment of obesity and CNS disorders)

RN 668476-24-2 CAPLUS

CN Pyrazolo[4,3-g][3]benzazepine, 4-chloro-3,6,7,8,9,10-hexahydro-3,8-dimethyl-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

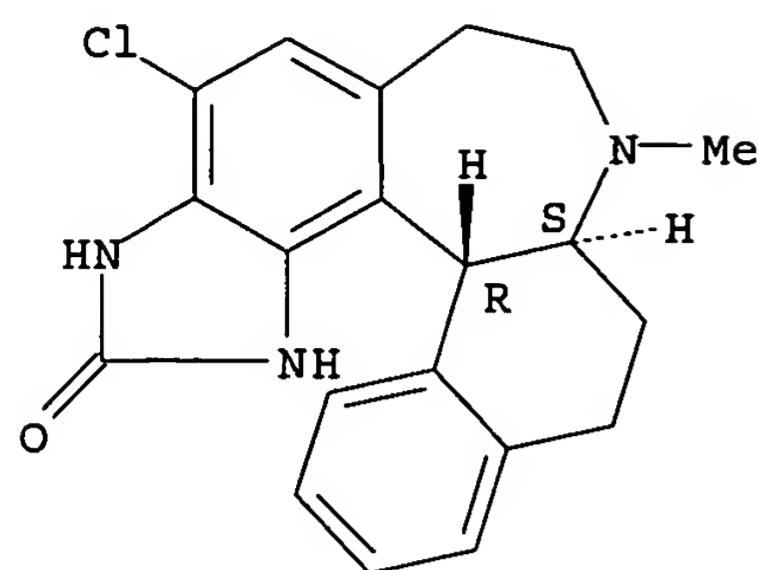
Absolute stereochemistry.



RN 668476-37-7 CAPLUS

CN Benzimidazo[5,4-d]benzo[g][1]benzazepin-2(1H)-one, 4-chloro-3,6,7,8,8a,9,10,14b-octahydro-8-methyl-, (8aS,14bR)- (9CI) (CA INDEX NAME)

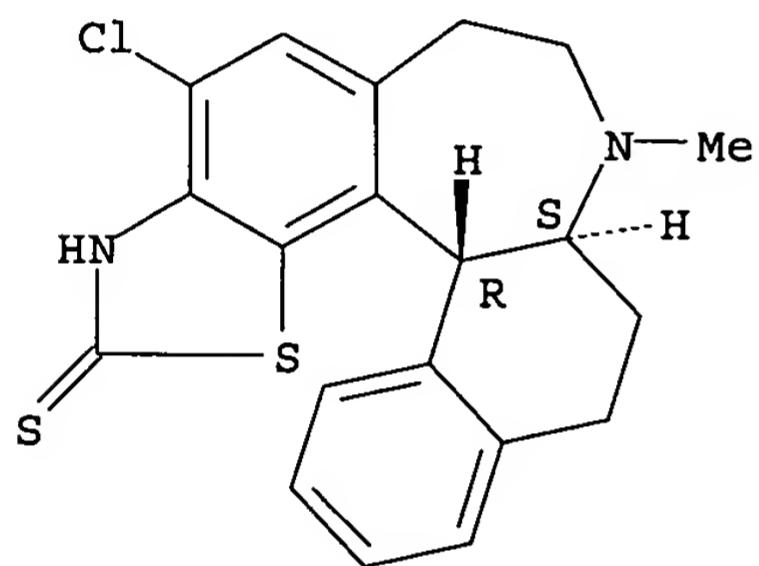
Absolute stereochemistry.



RN 668476-44-6 CAPLUS

CN 2H-Benz[g]benzothiazolo[6,7-d][1]benzazepine-2-thione,
4-chloro-3,6,7,8,8a,9,10,14b-octahydro-8-methyl-, (8aS,14bR)- (9CI) (CA
INDEX NAME)

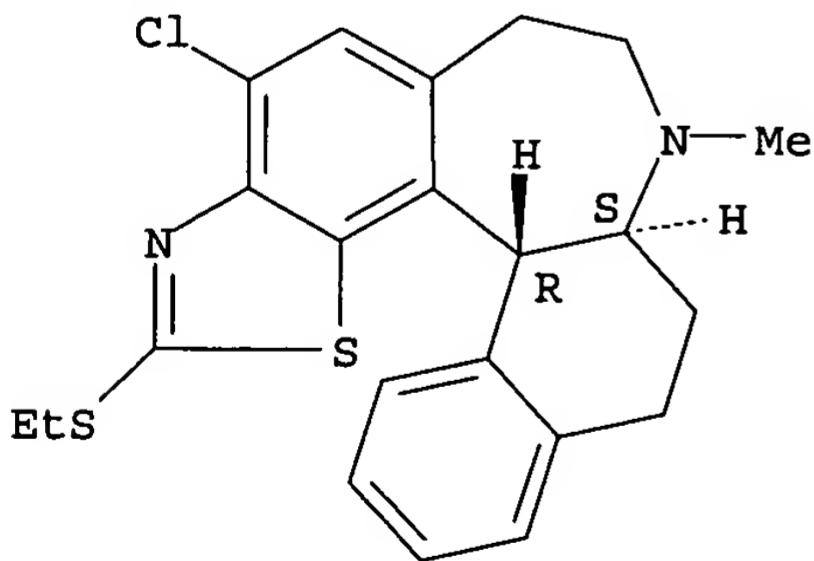
Absolute stereochemistry.



RN 668476-45-7 CAPLUS

CN 6H-Benz[g]benzothiazolo[6,7-d][1]benzazepine, 4-chloro-2-(ethylthio)-
7,8,8a,9,10,14b-hexahydro-8-methyl-, (8aS,14bR)- (9CI) (CA INDEX NAME)

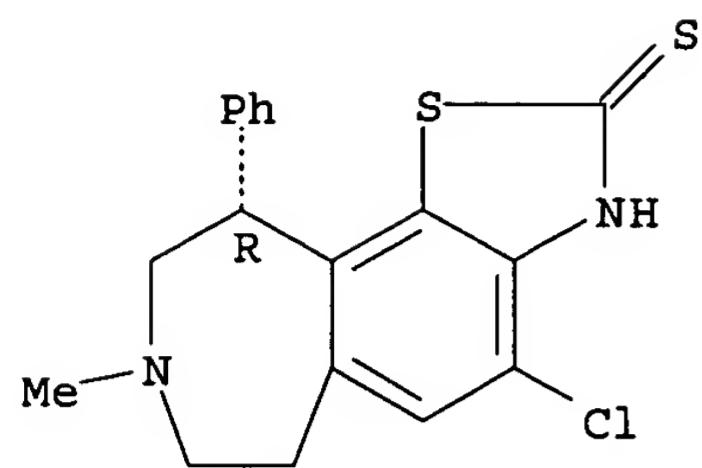
Absolute stereochemistry.



RN 668476-55-9 CAPLUS

CN 2H-Thiazolo[5,4-g][3]benzazepine-2-thione, 4-chloro-3,6,7,8,9,10-hexahydro-
8-methyl-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

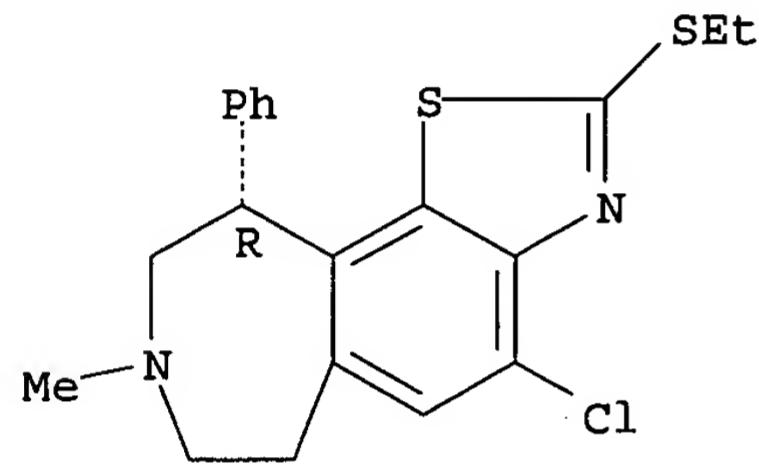
Absolute stereochemistry.



RN 668476-56-0 CAPLUS

CN 6H-Thiazolo[5,4-g][3]benzazepine, 4-chloro-2-(ethylthio)-7,8,9,10-tetrahydro-8-methyl-10-phenyl-, (10R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



REFERENCE COUNT:

3

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